LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010523\
Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : PE-9

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131931.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.369	34	48	57	rVB	153590	236940	13.83%	1.326%
2	4.998	490	495	506	rVB	502252	610335	35.64%	3.417%
3	5.410	561	565	573	rBV	1494170	1410620	82.36%	7.897%
4	6.440	735	740	743	rBV	1547189	1443427	84.28%	8.081%
5	6.798	798	801	805	rBV	449411	384938	22.48%	2.155%
6	7.063	843	846	853	rBV2	20189	21111	1.23%	0.118%
7	7.222	869	873	879	rVB3	20879	31144	1.82%	0.174%
8	7.375	894	899	901	rBV	1024312	986697	57.61%	5.524%
9	8.092	1018	1021	1023	rBV	557352	437864	25.57%	2.451%
10	8.110	1023	1024	1029	rVB	43341	36640	2.14%	0.205%
11	8.681	1118	1121	1125	rBV	28467	22415	1.31%	0.125%
12	9.175	1200	1205	1208	rBV	2131488	1712656	100.00%	9.588%
13	9.251	1212	1218	1220	rBV2	27884	31493	1.84%	0.176%
14	9.322	1227	1230	1234	rVB3	23271	23226	1.36%	0.130%
15	9.716	1293	1297	1300	rBV	19601	20450	1.19%	0.114%
16	9.851	1317	1320	1324	rVB	560941	493075	28.79%	2.760%
17	9.886	1324	1326	1329	rVB	34236	25304	1.48%	0.142%
18	10.057	1351	1355	1359	rBV5	21357	28652	1.67%	0.160%
19	10.098	1359	1362	1368	rVB3	21110	22481	1.31%	0.126%
20	10.575	1439	1443	1446	rVB	281122	253005	14.77%	1.416%
21	10.645	1451	1455	1459	rBV	967980	869034	50.74%	4.865%
22	10.769	1472	1476	1483	rVB4	48158	66855	3.90%	0.374%
23	11.080	1524	1529	1531	rBV3	33596	41158	2.40%	0.230%
24	11.192	1545	1548	1554	rBV3	31638	52378	3.06%	0.293%
25	11.345	1571	1574	1577	rBV	485867	432231	25.24%	2.420%
26	11.369	1577	1578	1581	rVV	100313	68032	3.97%	0.381%
27	11.422	1585	1587	1590	rVB	56811	43305	2.53%	0.242%
28	11.451	1590	1592	1596	rBV4	20507	31940	1.86%	0.179%
29	11.586	1611	1615	1620	rBV4	22346	53801	3.14%	0.301%
30	11.645	1622	1625	1628	rVV2	30042	28282	1.65%	0.158%
31	11.851	1658	1660	1662	rBV	46606	40412	2.36%	0.226%
32	11.880	1662	1665	1669	rVB	147560	123291	7.20%	0.690%
33	11.927	1671	1673	1676	rVB	55567	42142	2.46%	0.236%
34	11.963	1676	1679	1682	rBV	137434	147891	8.64%	0.828%
35	12.151	1708	1711	1713	rBV	44812	38023	2.22%	0.213%
36	12.174	1713	1715	1719	rVV3	60789	55963	3.27%	0.313%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010523\
Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.404	1751	1754	1756	rBV	66053	72676	4.24%	0.407%
38	12.492	1766	1769	1774	rBV	106184	112454	6.57%	0.630%
39	12.574	1779	1783	1786	rVV	1307000	1190059	69.49%	6.662%
40	12.804	1816	1822	1825	rBV	1283779	1335349	77.97%	7.476%
41	12.945	1842	1846	1849	rVB	1354161	1110782	64.86%	6.219%
42	13.233	1892	1895	1898	rBV2	129511	144570	8.44%	0.809%
43	13.857	1998	2001	2005	rVB	196562	209063	12.21%	1.170%
44	14.004	2022	2026	2029	rBV2	709458	1017719	59.42%	5.698%
45	14.039	2029	2032	2034	rVB	489736	449935	26.27%	2.519%
46	15.063	2203	2206	2210	rBV	563163	870453	50.82%	4.873%
47	15.374	2257	2259	2263	rBV	234437	265727	15.52%	1.488%
48	15.445	2267	2271	2276	rVB	531585	716464	41.83%	4.011%

Sum of corrected areas: 17862462

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Instrument :
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ClientSampleId :
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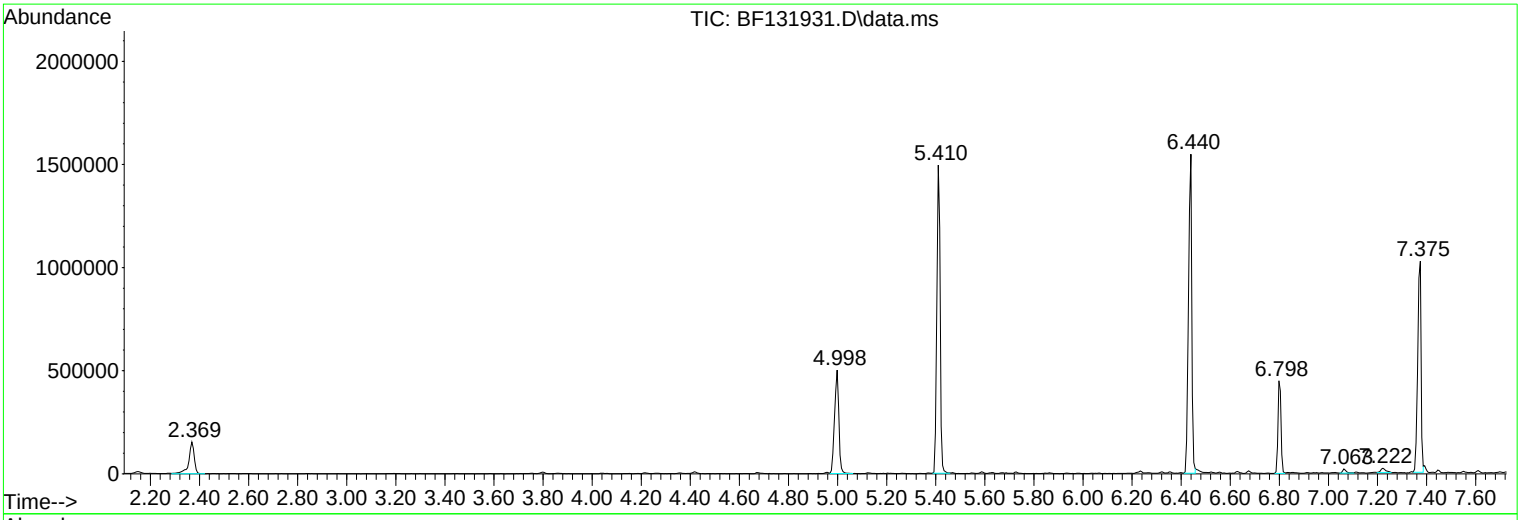
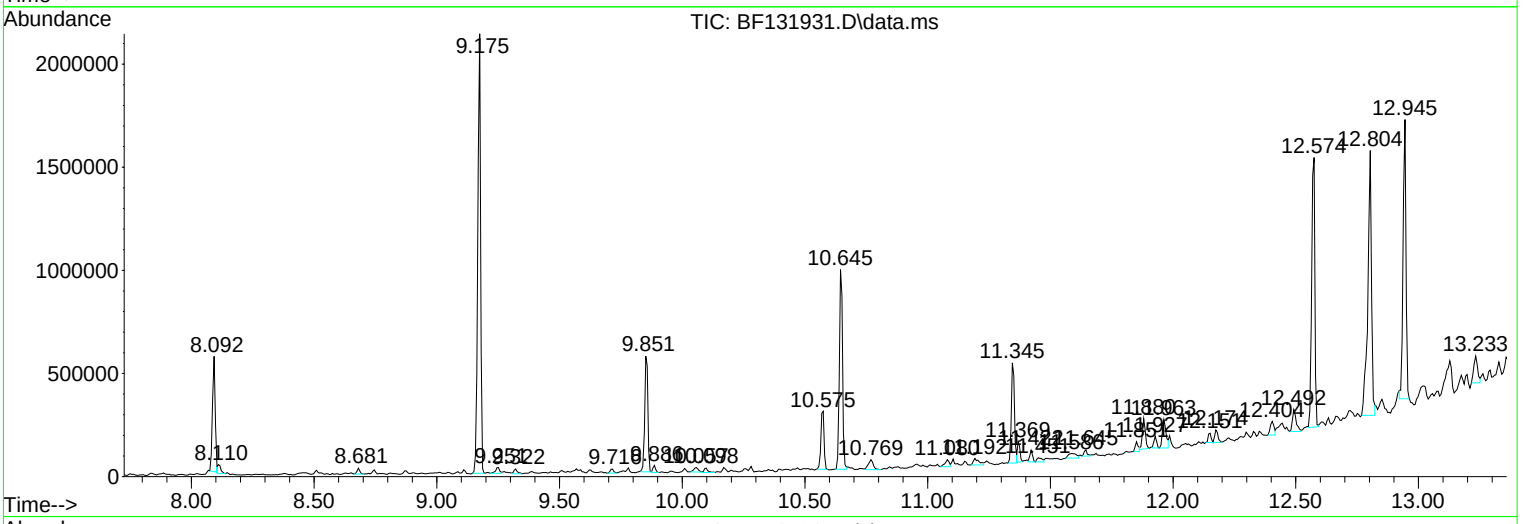
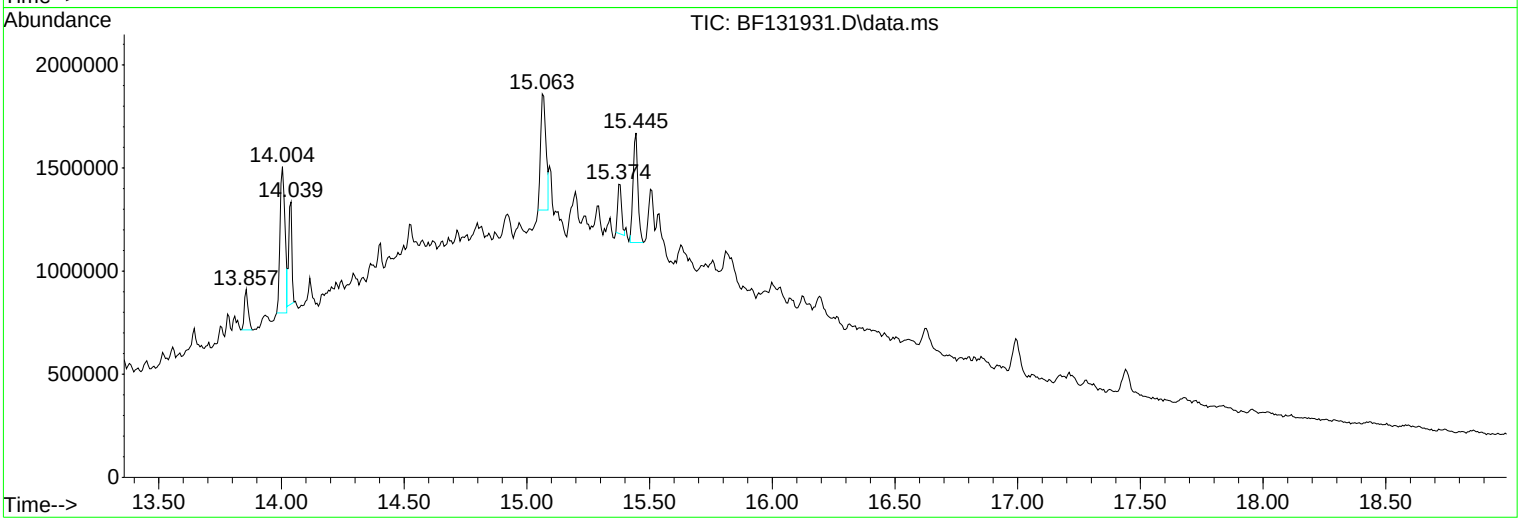
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P







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Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

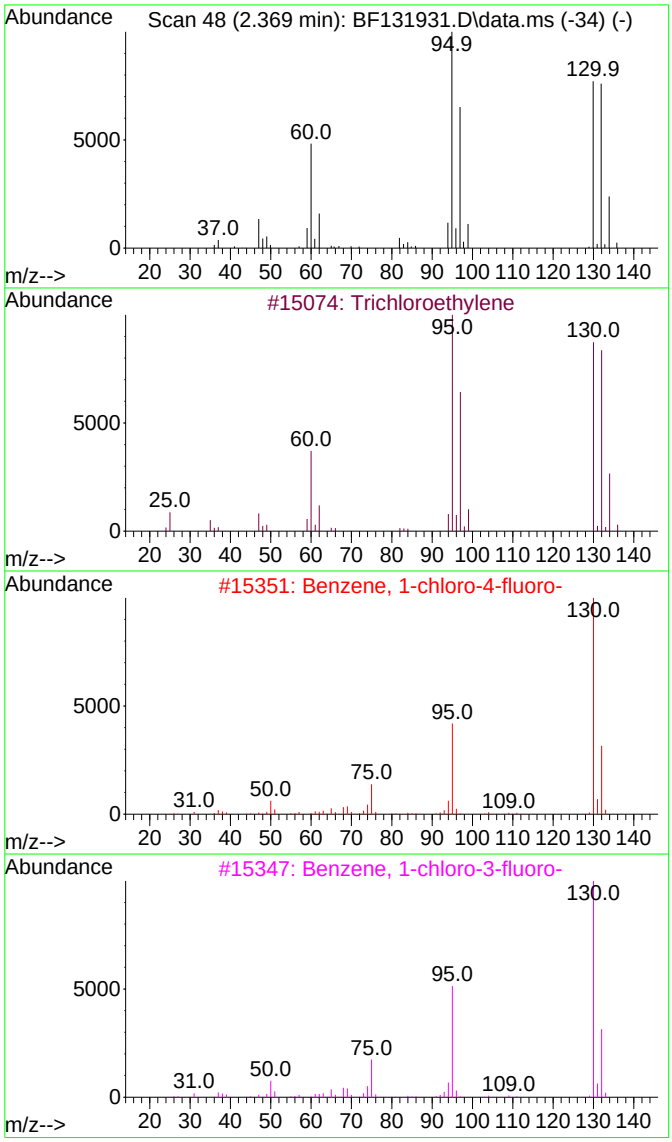
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

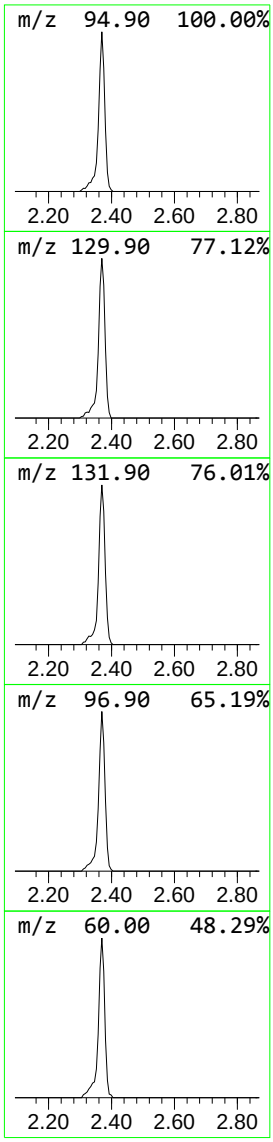
****

Peak Number 1 Trichloroethylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.369	12.31 ng	236940	1,4-Dichlorobenzene-d4	6.798

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trichloroethylene	130	C2HCl3	000079-01-6	97
2	Benzene, 1-chloro-4-fluoro-	130	C6H4ClF	000352-33-0	35
3	Benzene, 1-chloro-3-fluoro-	130	C6H4ClF	000625-98-9	35
4	Benzene, 1-chloro-2-fluoro-	130	C6H4ClF	000348-51-6	35
5	2-Fluoro-5-chloropyrimidine	132	C4H2ClFN2	062802-37-3	25





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Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

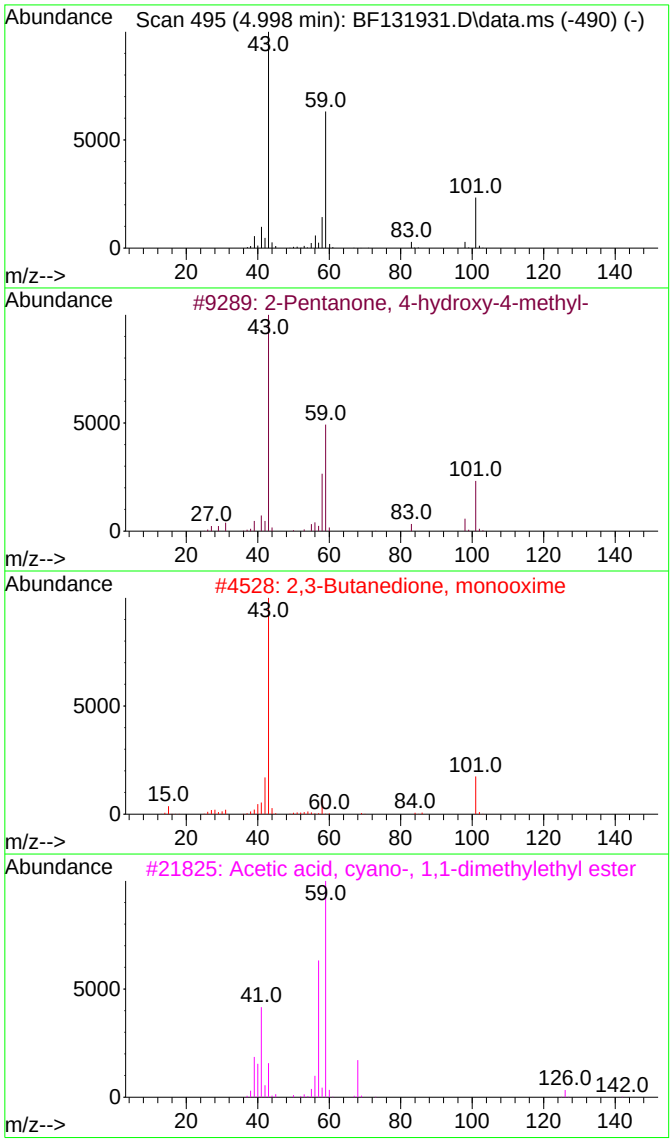
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

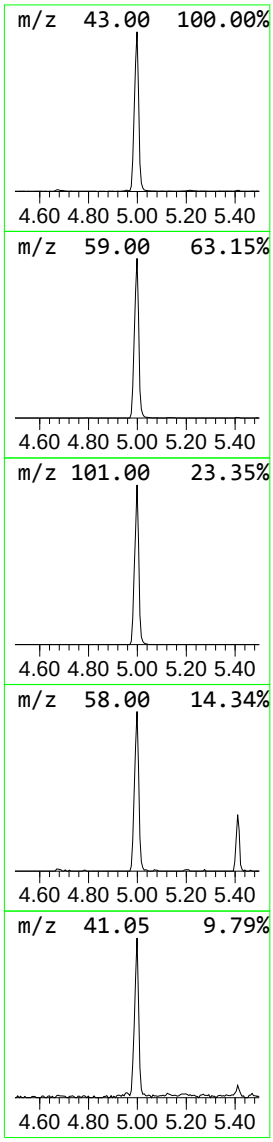
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.998	31.71 ng	610335	1,4-Dichlorobenzene-d4	6.798

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





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Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

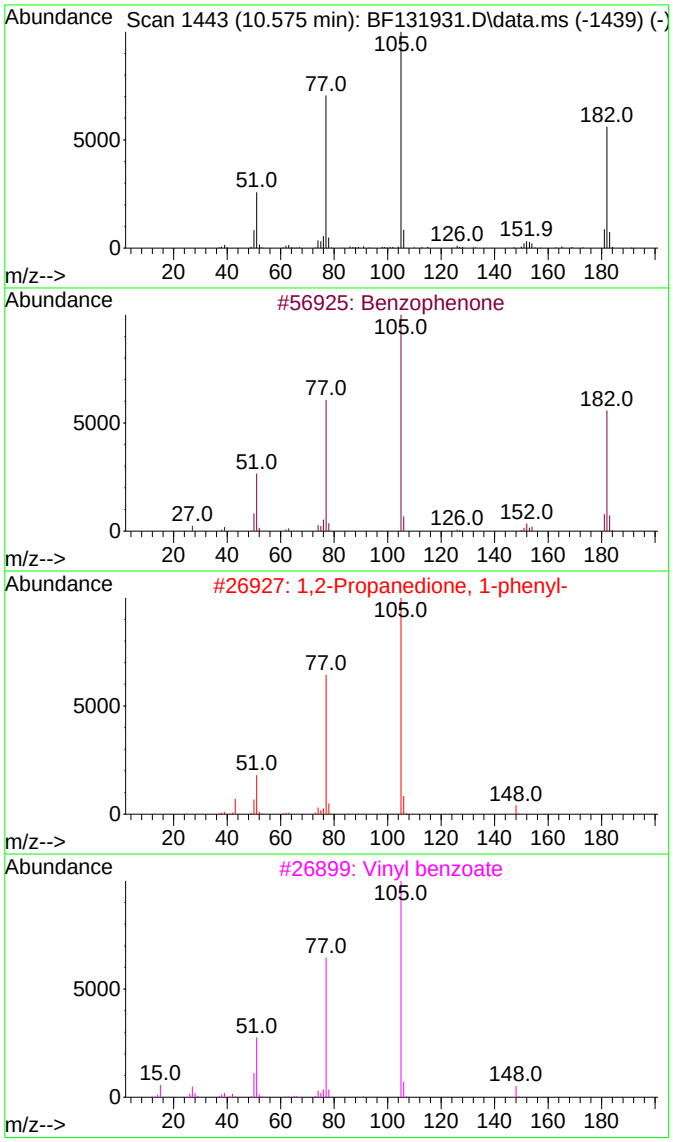
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

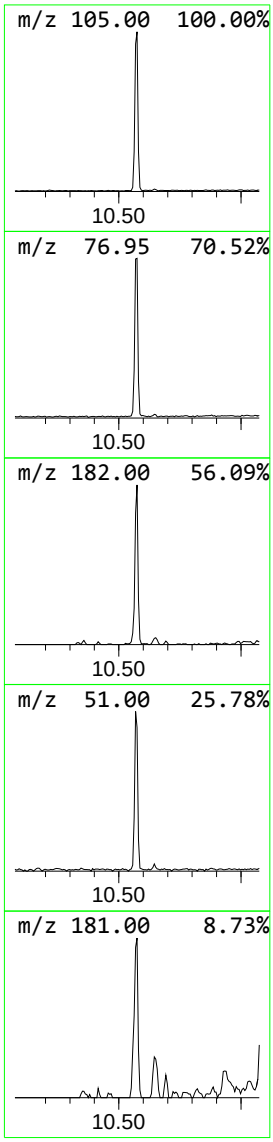
****

Peak Number 3 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.575	10.26 ng	253005	Acenaphthene-d10	9.851

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	47
3	Vinyl benzoate	148	C9H8O2	000769-78-8	47
4	Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	47
5	Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010523\
Data File : BF131931.D
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Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

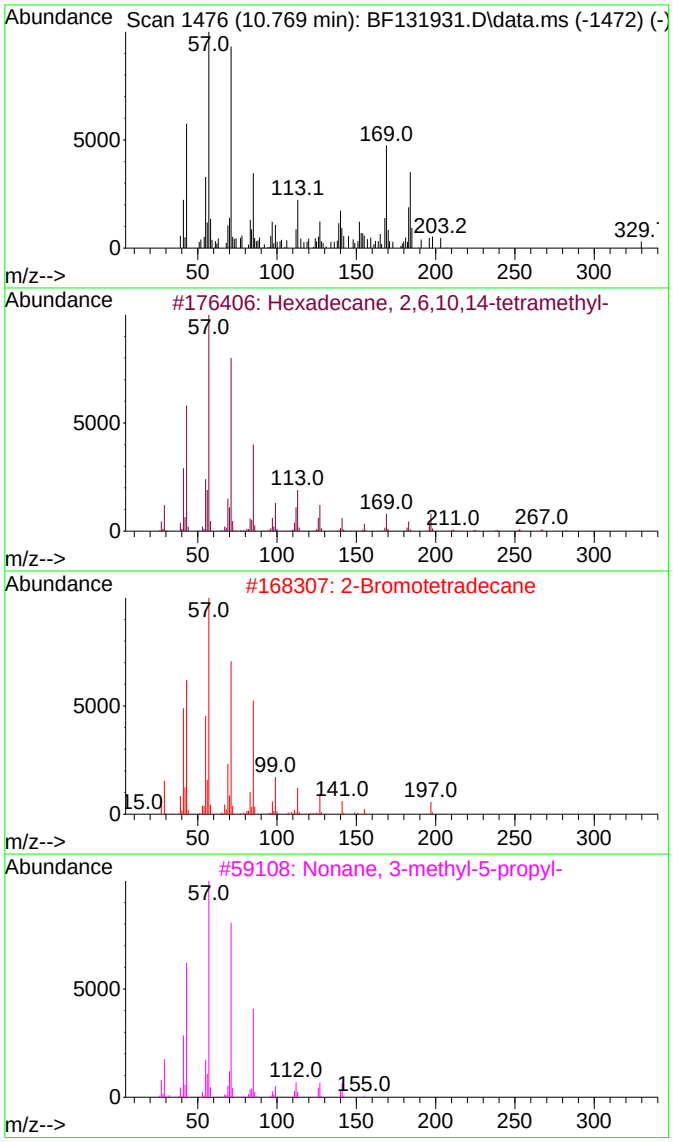
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

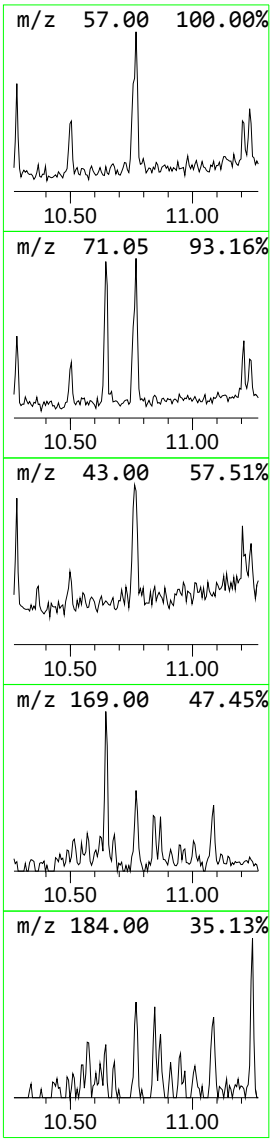
****

Peak Number 4 unknown10.769 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.769	3.09 ng	66855	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	49
2	2-Bromotetradecane	276	C14H29Br	074036-95-6	38
3	Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	38
4	Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	38
5	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010523\
Data File : BF131931.D
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Operator : CG\JU
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

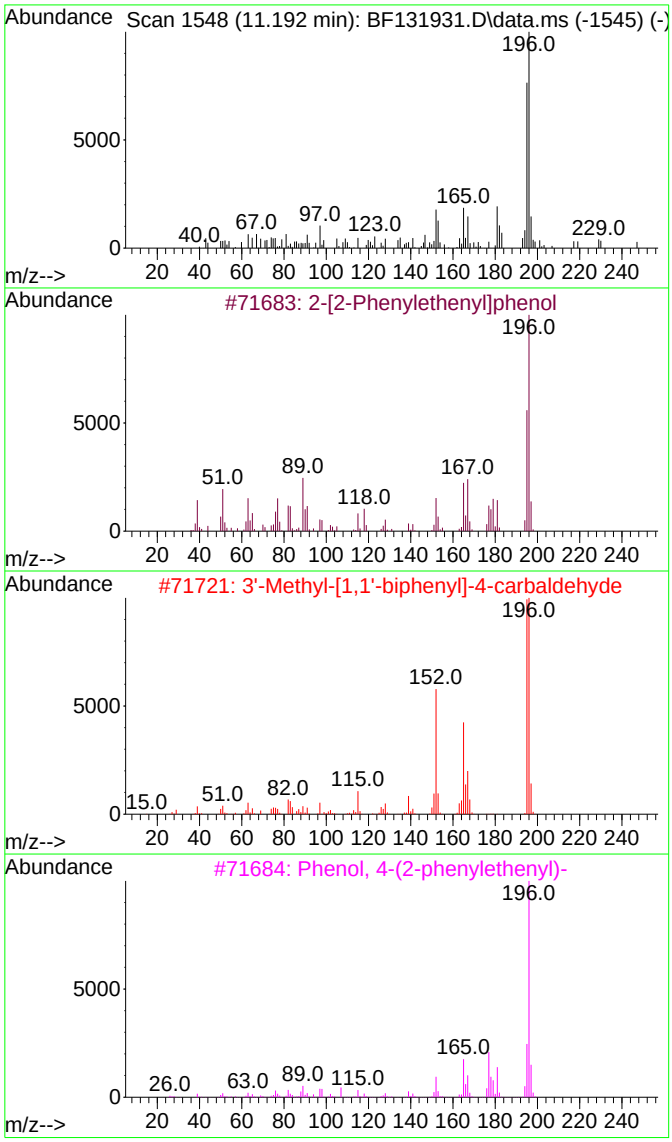
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

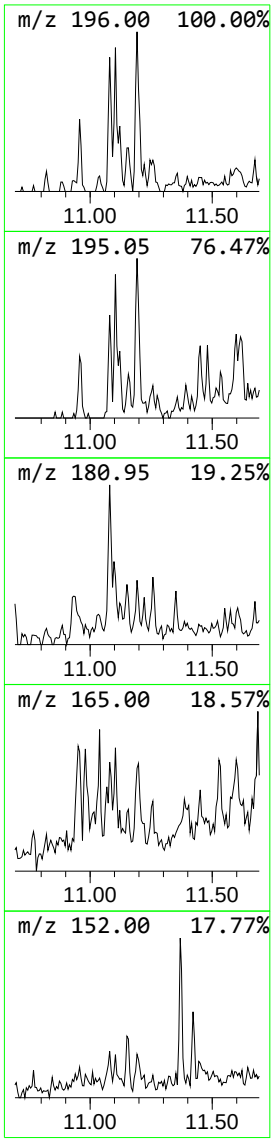
****

Peak Number 5 2-[2-Phenylethenyl]phenol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.192	2.42 ng	52378	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-[2-Phenylethenyl]phenol	196	C14H12O	042224-48-6	80
2	3'-Methyl-[1,1'-biphenyl]-4-carb...	196	C14H12O	400744-83-4	68
3	Phenol, 4-(2-phenylethenyl)-	196	C14H12O	003839-46-1	52
4	Naphtho[2,1-b]furan, 1,2-dimethyl-	196	C14H12O	129812-23-3	49
5	8-Dimethylaminonaphthalene-1-car...	196	C13H12N2	128644-69-9	47





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Data File : BF131931.D
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Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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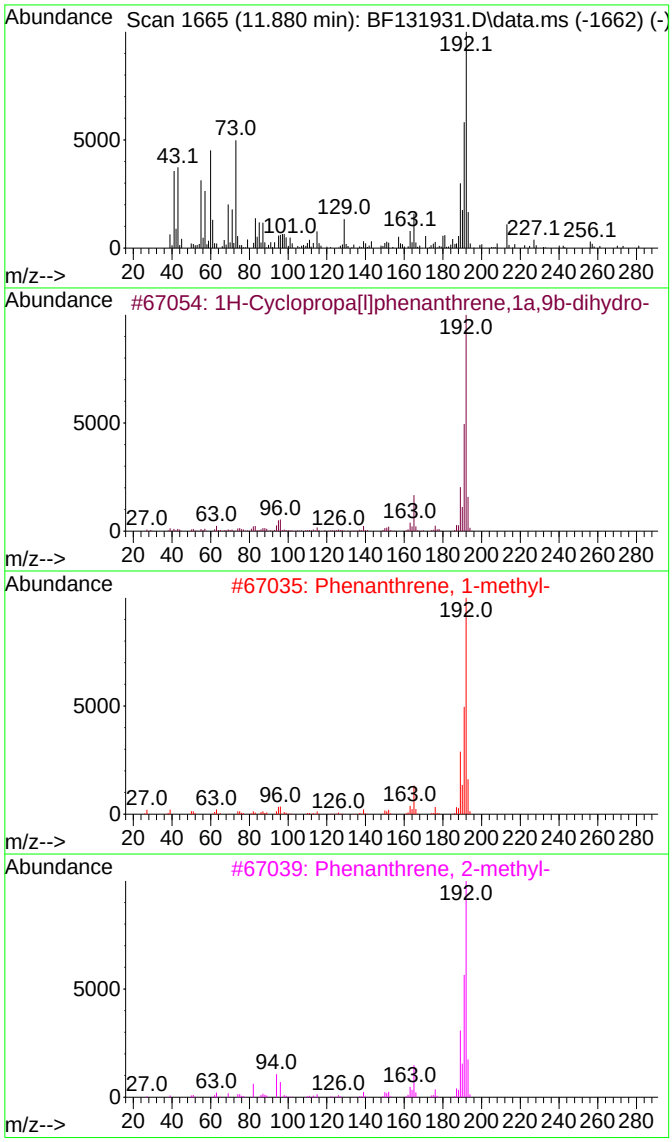
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

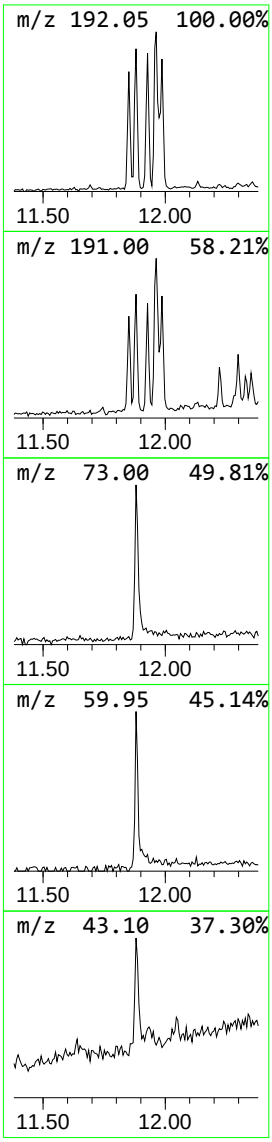
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Cyclopropa[l]phenanthren... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	5.70 ng	123291	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	90
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	78





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Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

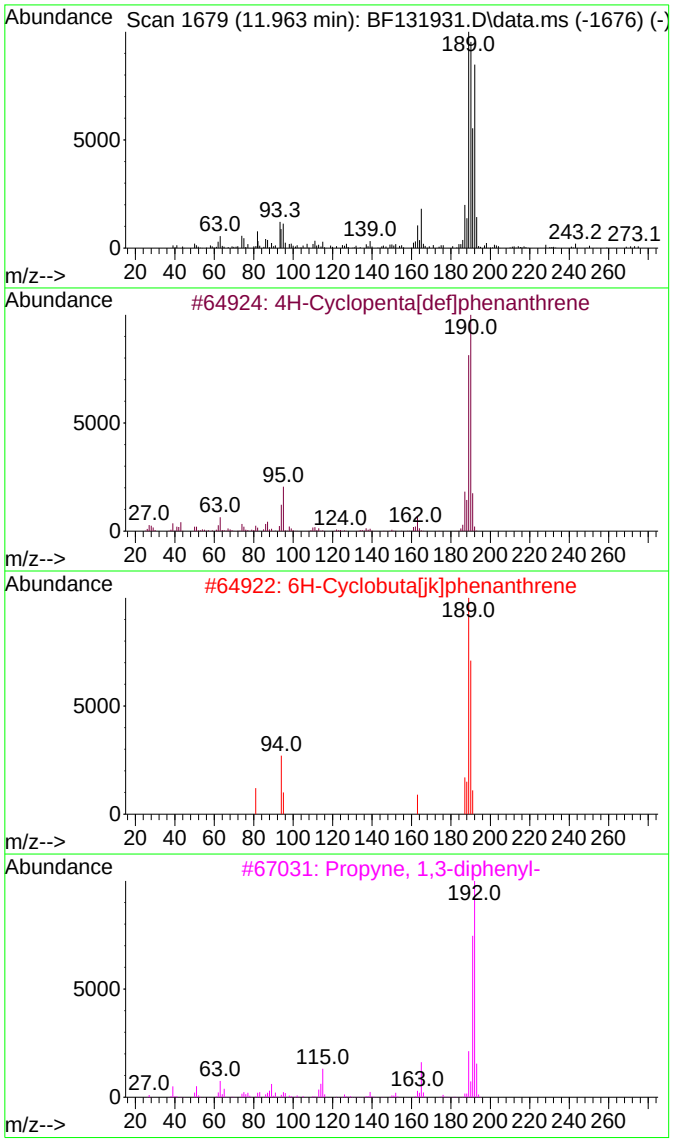
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

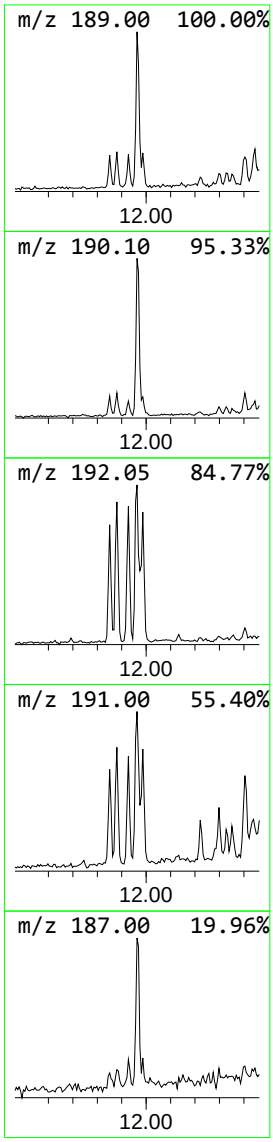
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.963	6.84 ng	147891	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
3	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	46
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	42
5	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	42





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010523\
Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

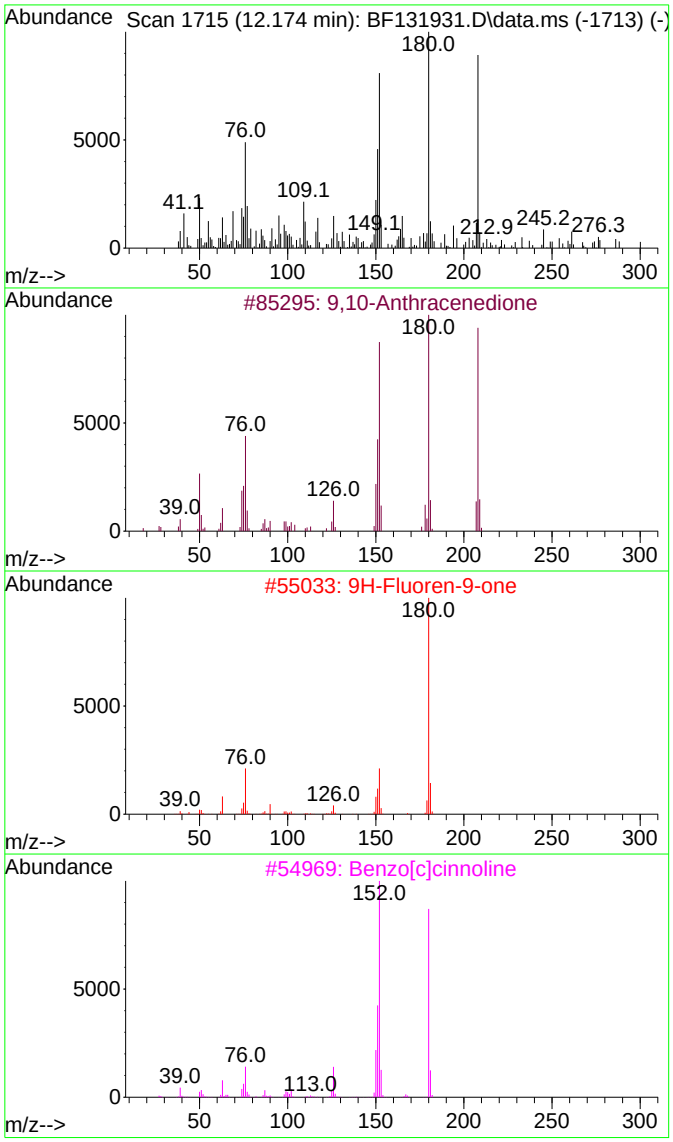
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

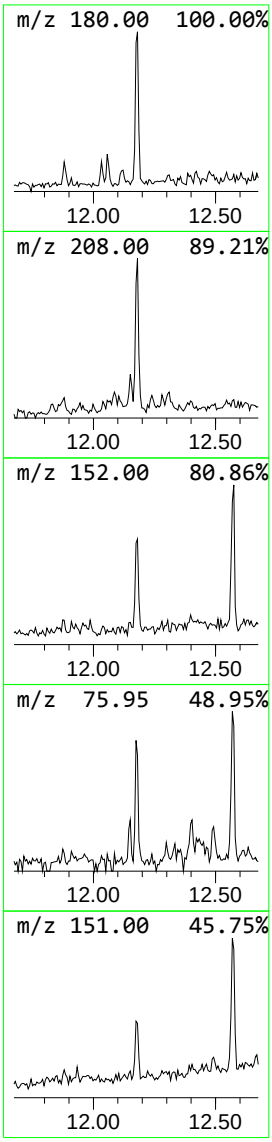
****

Peak Number 9 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.175	2.59 ng	55963	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
2	9H-Fluoren-9-one	180	C13H8O	000486-25-9	91
3	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	60
4	Benzo[h]cinnoline	180	C12H8N2	000230-31-9	50
5	9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	49





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Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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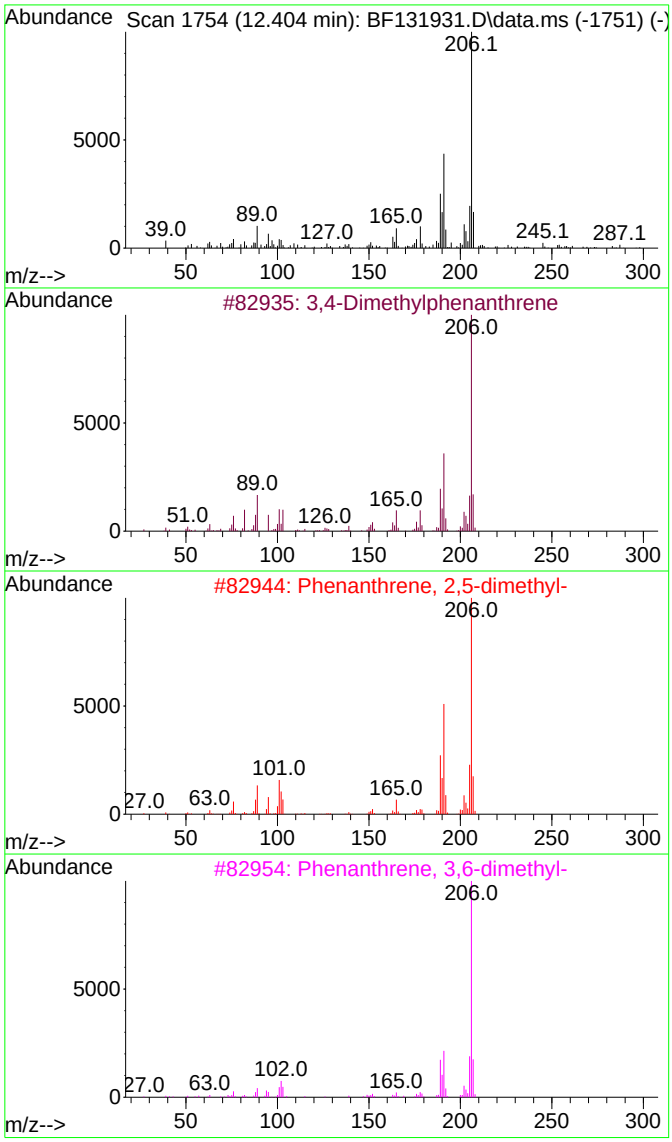
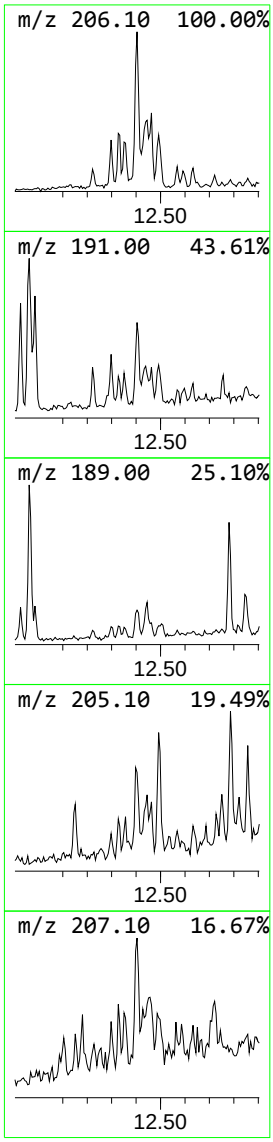
Peak Number 10 3,4-Dimethylphenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.404	3.36 ng	72676	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	94
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91
5	di-p-Tolylacetylene	206	C16H14	002789-88-0	90





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Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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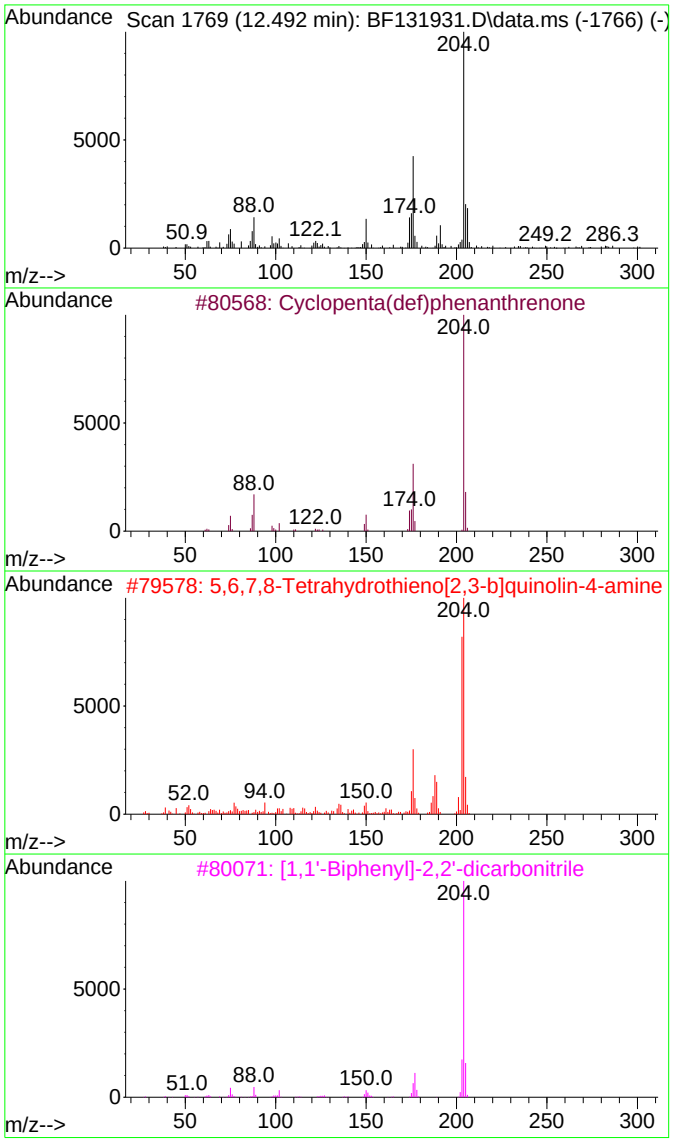
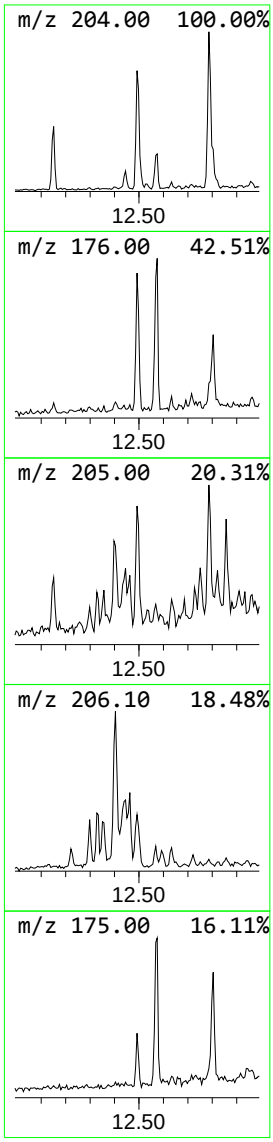
Peak Number 11 Cyclopenta(def)phenanthrenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.492	5.20 ng	112454	Phenanthrene-d10	11.345

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	81
2	5,6,7,8-Tetrahydrothieno[2,3-b]q...	204	C11H12N2S	122914-50-5	47
3	[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	46
4	[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	43
5	Methyl 3,4,5-trihydroxycyclohexe...	404	C17H36O5Si3	1000493-58-0	43





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Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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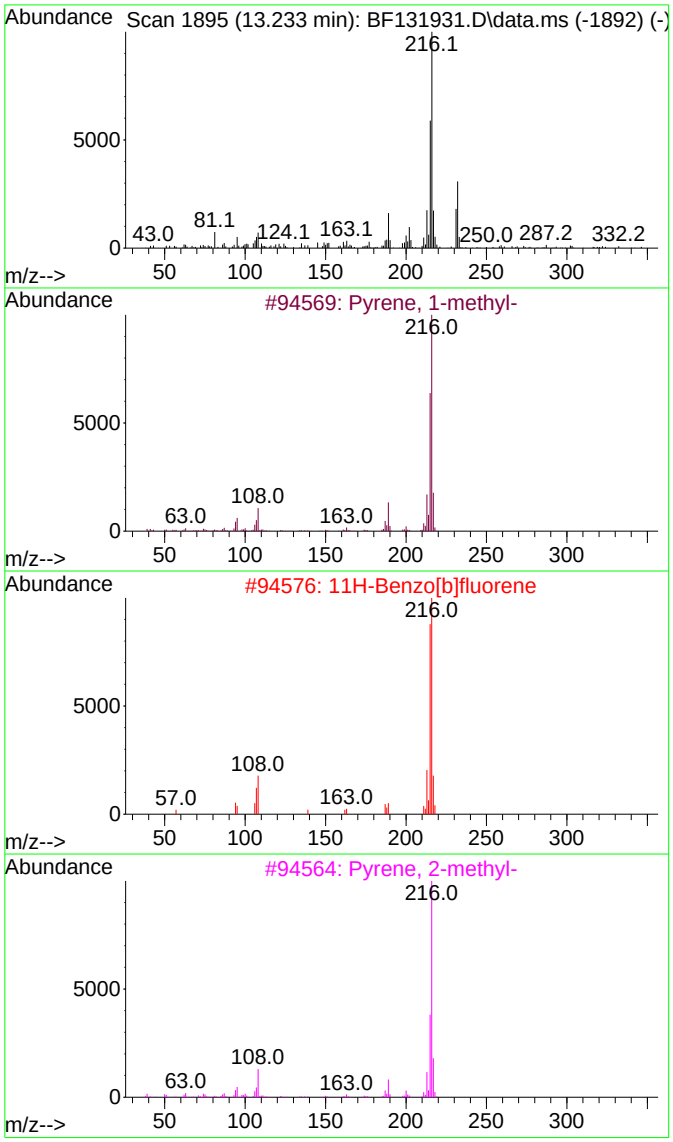
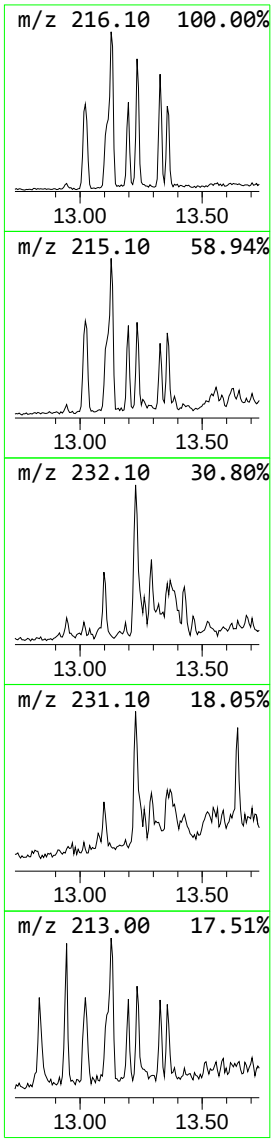
Peak Number 12 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.233	2.84 ng	144570	Chrysene-d12	14.010

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	89
4	Pyrene, 4-methyl-	216	C17H12	003353-12-6	83
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	70





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Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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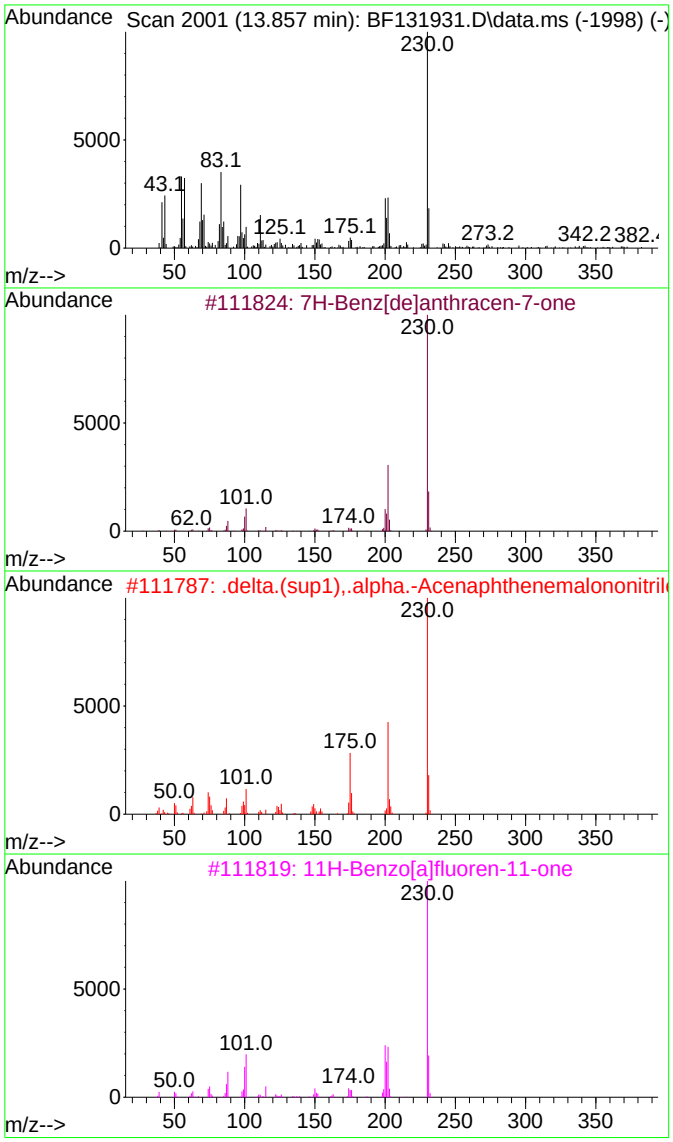
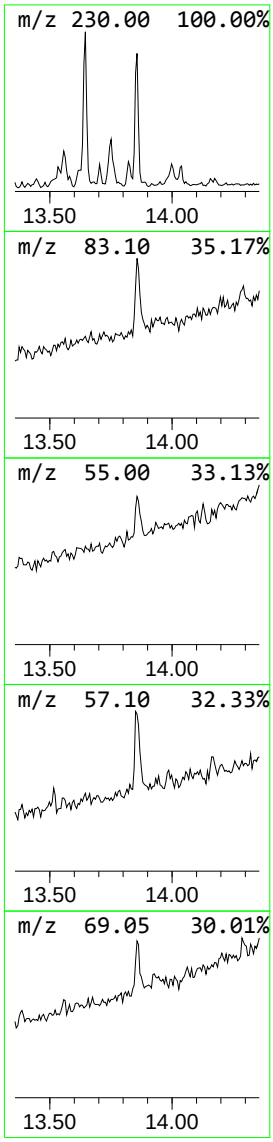
Peak Number 13 7H-Benz[de]anthracen-7-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	4.11 ng	209063	Chrysene-d12	14.010

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	64
2	.delta.(sup1),.alpha.-Acenaphthe...	230	C15H6N2O	014619-86-4	53
3	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	52
4	benzonitrile, 2-(3-isoquinolinyl)-	230	C16H10N2	1000401-00-3	46
5	1H-isoindole, 1-(1H-isoindol-1-y...	230	C16H10N2	1000396-37-4	43





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Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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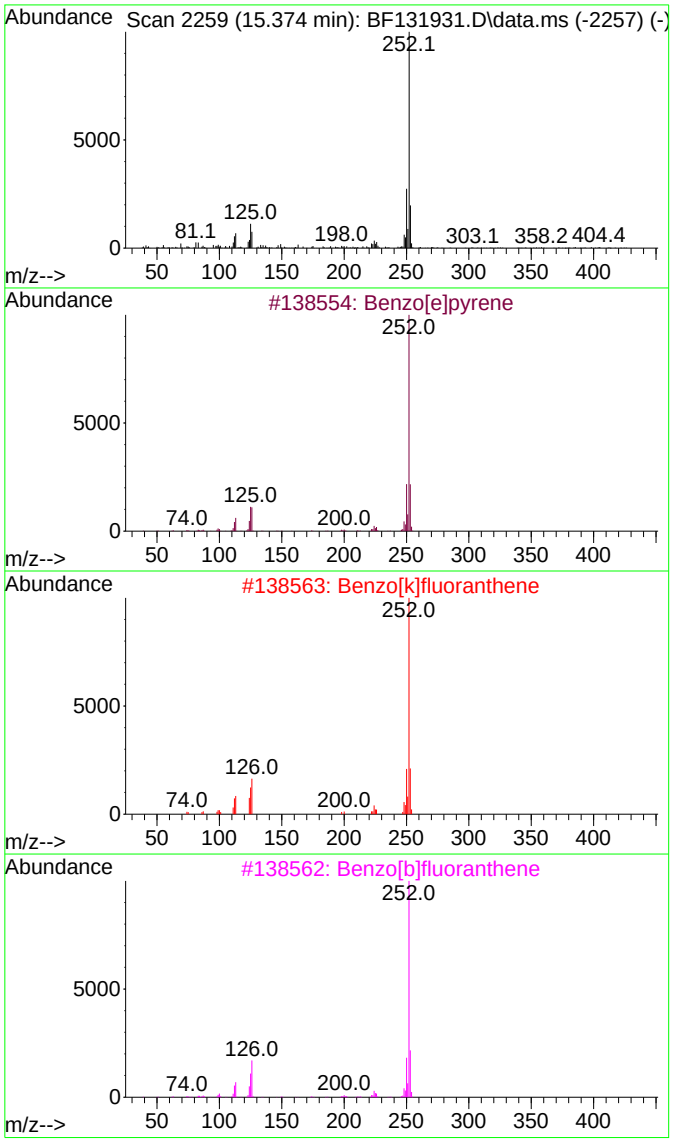
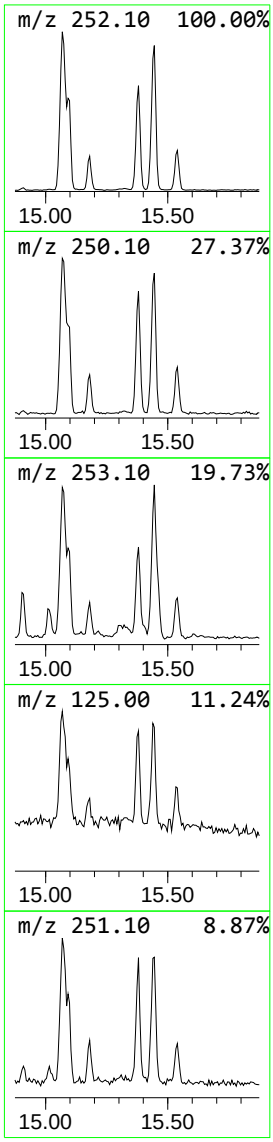
Peak Number 14 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.374	7.42 ng	265727	Perylene-d12	15.504

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3	Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4	Benzo[a]pyrene	252	C20H12	000050-32-8	90
5	Benzo[j]fluoranthene	252	C20H12	000205-82-3	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010523\
Data File : BF131931.D
Acq On : 05 Jan 2023 23:47
Operator : CG\JU
Sample : N6243-09
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PE-9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Trichloroethylene	2.369	12.3	ng	236940	1	6.798	384938	20.0
2-Pentanone, 4-...	4.998	31.7	ng	610335	1	6.798	384938	20.0
Benzophenone	10.575	10.3	ng	253005	3	9.851	493075	20.0
unknown10.769	10.769	3.1	ng	66855	4	11.345	432231	20.0
2-[2-Phenylethe...	11.192	2.4	ng	52378	4	11.345	432231	20.0
1H-Cyclopropa[l...	11.880	5.7	ng	123291	4	11.345	432231	20.0
4H-Cyclopenta[d...	11.963	6.8	ng	147891	4	11.345	432231	20.0
9,10-Anthracene...	12.175	2.6	ng	55963	4	11.345	432231	20.0
3,4-Dimethylphe...	12.404	3.4	ng	72676	4	11.345	432231	20.0
Cyclopenta(def)...	12.492	5.2	ng	112454	4	11.345	432231	20.0
Pyrene, 1-methyl-	13.233	2.8	ng	144570	5	14.010	1017720	20.0
7H-Benz[de]anth...	13.857	4.1	ng	209063	5	14.010	1017720	20.0
Benzo[e]pyrene	15.374	7.4	ng	265727	6	15.504	716464	20.0